NEWS 30

JAN 16

Connecting via Winsock to STN

```
Welcome to STN International! Enter x:x
LOGINID: SSPTAEAL1624
PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2
* * * * * * * * * *
                     Welcome to STN International
                                                    * * * * * * * * * *
NEWS
                 Web Page for STN Seminar Schedule - N. America
NEWS
      2 AUG 06 CAS REGISTRY enhanced with new experimental property tags
NEWS
      3
         AUG 06
                 FSTA enhanced with new thesaurus edition
NEWS
         AUG 13
                 CA/CAplus enhanced with additional kind codes for granted
                 patents
NEWS
         AUG 20
                 CA/CAplus enhanced with CAS indexing in pre-1907 records
NEWS
         AUG 27
                 Full-text patent databases enhanced with predefined
                 patent family display formats from INPADOCDB
         AUG 27
                 USPATOLD now available on STN
NEWS
NEWS 8
         AUG 28
                 CAS REGISTRY enhanced with additional experimental
                 spectral property data
NEWS 9
         SEP 07
                 STN AnaVist, Version 2.0, now available with Derwent
                 World Patents Index
NEWS 10 SEP 13
                 FORIS renamed to SOFIS
NEWS 11 SEP 13
                 INPADOCDB enhanced with monthly SDI frequency
NEWS 12 SEP 17
                 CA/CAplus enhanced with printed CA page images from
                 1967-1998
NEWS 13
         SEP 17
                 CAplus coverage extended to include traditional medicine
NEWS 14 SEP 24
                 EMBASE, EMBAL, and LEMBASE reloaded with enhancements
NEWS 15 OCT 02 CA/Caplus enhanced with pre-1907 records from Chemisches
                 Zentralblatt
NEWS 16 OCT 19
                 BEILSTEIN updated with new compounds
NEWS 17 NOV 15 Derwent Indian patent publication number format enhanced
NEWS 18 NOV 19 WPIX enhanced with XML display format
NEWS 19 NOV 30 ICSD reloaded with enhancements
NEWS 20 DEC 04 LINPADOCDB now available on STN
NEWS 21 DEC 14 BEILSTEIN pricing structure to change
NEWS 22 DEC 17 USPATOLD added to additional database clusters
NEWS 23
         DEC 17
                 IMSDRUGCONF removed from database clusters and STN
NEWS 24
         DEC 17
                 DGENE now includes more than 10 million sequences
NEWS 25
         DEC 17
                 TOXCENTER enhanced with 2008 MeSH vocabulary in
                 MEDLINE segment
         DEC 17
NEWS 26
                 MEDLINE and LMEDLINE updated with 2008 MeSH vocabulary
NEWS 27
         DEC 17
                 CA/CAplus enhanced with new custom IPC display formats
NEWS 28
         DEC 17
                 STN Viewer enhanced with full-text patent content
                  from USPATOLD
NEWS 29
                 STN pricing information for 2008 now available
         JAN 02
```

CAS patent coverage enhanced to include exemplified

<12/04/2007> Erich Leese

prophetic substances

NEWS EXPRESS 19 SEPTEMBER 2007: CURRENT WINDOWS VERSION IS V8.2, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 19 SEPTEMBER 2007.

NEWS HOURS STN Operating Hours Plus Help Desk Availability

NEWS LOGIN Welcome Banner and News Items

NEWS IPC8 For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

FILE 'HOME' ENTERED AT 15:01:03 ON 18 JAN 2008

=> file reg COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 15:01:15 ON 18 JAN 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 17 JAN 2008 HIGHEST RN 1000264-70-9 DICTIONARY FILE UPDATES: 17 JAN 2008 HIGHEST RN 1000264-70-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

Please note that search-term pricing does apply when conducting ${\tt SmartSELECT}$ searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

Uploading C:\Program Files\Stnexp\Queries\10596851.str

G1:Cb, Ak

G2:H,Cb,Ak

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:CLASS 20:CLASS 21:CLASS 22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS 31:CLASS 32:CLASS 34:CLASS

L1 STRUCTURE UPLOADED

=> s 11 full

FULL SEARCH INITIATED 15:02:15 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 203 TO ITERATE

100.0% PROCESSED 203 ITERATIONS

90 ANSWERS

SEARCH TIME: 00.00.01

L2 90 SEA SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 178.82 179.03

FILE 'CAPLUS' ENTERED AT 15:02:19 ON 18 JAN 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 18 Jan 2008 VOL 148 ISS 4 FILE LAST UPDATED: 17 Jan 2008 (20080117/ED)

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

http://www.cas.org/infopolicy.html

=> s 12 full L3 2 L2

=> d ibib abs hitstr tot

L3 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2005:638860 CAPLUS

DOCUMENT NUMBER: 143:153402

TITLE: Preparation of diarylmethylpiperazines as δ receptor ligands for the treatment of pain

INVENTOR(S): Brown, William; Griffin, Andrew

PATENT ASSIGNEE(S): Astrazeneca AB, Swed. SOURCE: PCT Int. Appl., 69 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PA:	TENT	NO.			KIND		DATE			APPI	LICAT	DATE						
WO	2005					20050721			 WO 2	2005-	20050105							
											BG,							
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,	
		GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KP,	KR,	KZ,	LC,	
		LK,	LR,	LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	NA,	NI,	
		NO,	NZ,	OM,	PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,	SG,	SK,	SL,	SY,	
		ТJ,	TM,	TN,	TR,	TT,	TZ,	UA,	UG,	US,	UZ,	VC,	VN,	YU,	ZA,	ZM,	ZW	
	RW:	BW,	GH,	GM,	ΚE,	LS,	MW,	MΖ,	NA,	SD,	SL,	SZ,	TZ,	UG,	ZM,	ZW,	AM,	
		ΑZ,	BY,	KG,	KΖ,	MD,	RU,	ТJ,	TM,	ΑT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,	
		EE,	ES,	FI,	FR,	GB,	GR,	ΗU,	ΙE,	IS,	IT,	LT,	LU,	MC,	NL,	PL,	PT,	
		RO,	SE,	SI,	SK,	TR,	BF,	ВJ,	CF,	CG,	CI,	CM,	GA,	GN,	GQ,	G₩,	ML,	
					TD,													
AU	2005	2040	10							AU 2	2005-		20050105					
					A1 20050721													
EP								EP 2005-704688										
	R:	ΑT,	BE,	CH,	DE,	DK,	ES,	FR,	GB,	GR,	IT,	LI,	LU,	NL,	SE,	MC,	PT,	
											CZ,							
CN	CN 1926122 BR 2005006702 JP 2007517873						2007	0307		CN 2	2005-		20050105					
BR	BR 2005006702						2007	0502		BR 2	2005-		20050105					
JP	JP 2007517873						2007	0705		JP 2	2006-		20050105					
	IN 2006DN03738												20060629					
	MX 2006PA07664														20060703			
	NO 2006003619														20060809			
	US 2007293502						2007	1220		US 2	2007-	5968	51		2	0070	529	
PRIORIT	IORITY APPLN. INFO.:																	
							WO 2005-SE14								W 20050105			
OTHER SO						CASREACT 143:153402; MARPAT 143:153402												

^{*} STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB Title compds. I [R1 = alkyl, alkenyl, cycloalkyl, etc.; R2 = H, alkyl, cycloalkyl, etc.; R3 = alkyl, cycloalkyl] and their pharmaceutically acceptable salts were prepared For example, N-alkylation of piperazine II (R1 =H) with bromoethyl Me ether afforded the hCL salt of claimed diarylmethylpiperazine II (R1 = CH2CH2OCH3) in 68% yield. In human δ receptor assays, certain examples of compds. I exhibited IC50 values ranging from 0.2-3.7 nM, with an average of 1 nM (sic).

```
ΤТ
     859634-99-4P 859635-00-0P 859635-01-1P
     859635-02-2P 859635-03-3P 859635-04-4P
     859635-05-5P 859635-06-6P 859635-07-7P
     859635-08-8P 859635-09-9P 859635-10-2P
     859635-11-3P 859635-12-4P 859635-13-5P
     859635-14-6P 859635-15-7P 859635-16-8P
     859635-17-9P 859635-18-0P 859843-90-6P
     859843-91-7P 859843-92-8P 859843-93-9P
     859843-94-0P 859843-95-1P 859843-96-2P
     859843-97-3P 859843-98-4P 859843-99-5P
     859844-00-1P 859844-01-2P
     RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
     (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
     (Uses)
        (preparation of diarylmethylpiperazines as \delta receptor ligands for
        treatment of pain)
RN
     859634-99-4 CAPLUS
CN
     Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(2-
     methoxyethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester,
     monohydrochloride (9CI) (CA INDEX NAME)
```

Absolute stereochemistry. Rotation (+).

● HCl

```
RN 859635-00-0 CAPLUS
CN Carbamic acid, [3-[(S)-(4-butyl-1-piperazinyl)][4-
[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, methyl ester (9CI) (CA
INDEX NAME)
```

Absolute stereochemistry. Rotation (+).

RN 859635-01-1 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl](4-pentyl-1-piperazinyl)methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 859635-02-2 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl](4-propyl-1-piperazinyl)methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 859635-03-3 CAPLUS

CN Carbamic acid, [3-[(S)-[4-(cyclopropylmethyl)-1-piperazinyl]][4[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, methyl ester (9CI) (CA
INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 859635-04-4 CAPLUS

CN Carbamic acid, [3-[(S)-[4-(cyclobutylmethyl)-1-piperazinyl][4-[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 859635-05-5 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl]][4-(2-methoxyethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859635-06-6 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-ethoxyethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

HC1

RN 859635-07-7 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(3-methoxypropyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859635-08-8 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl](4-propyl-1-piperazinyl)methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859635-09-9 CAPLUS

CN Carbamic acid, [3-[(R)-(4-butyl-1-piperazinyl)][4[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, methyl ester (9CI) (CA
INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859635-10-2 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl](4-pentyl-1-piperazinyl)methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859635-11-3 CAPLUS

CN Carbamic acid, [3-[(R)-[4-(cyclopropylmethyl)-1-piperazinyl][4[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, methyl ester (9CI) (CF
INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859635-12-4 CAPLUS

CN Carbamic acid, [3-[(R)-[4-(cyclobutylmethyl)-1-piperazinyl][4[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, methyl ester (9CI) (CA
INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859635-13-5 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-methoxyethyl)-1-piperazinyl]methyl]phenyl]-, ethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859635-14-6 CAPLUS

CN Carbamic acid, [3-[(R)-(4-butyl-1-piperazinyl)][4[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, ethyl ester (9CI) (CA
INDEX NAME)

Absolute stereochemistry. Rotation (-).

$$\begin{array}{c|c} n-Bu \\ N \\ N \\ R \\ \end{array}$$

RN 859635-15-7 CAPLUS

CN Carbamic acid, [3-[(R)-[4-(cyclopropylmethyl)-1-piperazinyl]][4[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, ethyl ester (9CI) (CA
INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859635-16-8 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl](4-propyl-1-piperazinyl)methyl]phenyl]-, ethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859635-17-9 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl](4-ethyl-1-piperazinyl)methyl]phenyl]-, ethyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

$$\begin{array}{c|c} \text{Et} \\ N \\ N \\ \end{array}$$

● HCl

RN 859635-18-0 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl](4-methyl-1-piperazinyl)methyl]phenyl]-, ethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859843-90-6 CAPLUS

CN Carbamic acid, [3-[(S)-(4-butyl-1-piperazinyl)]4[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, methyl ester,
monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

● HCl

RN 859843-91-7 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl](4-pentyl-1-piperazinyl)methyl]phenyl]-, methyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Erich Leese

Absolute stereochemistry. Rotation (+).

<12/04/2007>

RN 859843-92-8 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-methoxyethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

● HCl

RN 859843-93-9 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl](4-propyl-1-piperazinyl)methyl]phenyl]-, methyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

● HCl

RN 859843-94-0 CAPLUS

CN Carbamic acid, [3-[(R)-(4-butyl-1-piperazinyl)][4-[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, methyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

● HCl

RN 859843-95-1 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl](4-pentyl-1-piperazinyl)methyl]phenyl]-, methyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859843-96-2 CAPLUS

CN Carbamic acid, [3-[(R)-[4-(cyclopropylmethyl)-1-piperazinyl][4-[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, methyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

● HCl

RN 859843-97-3 CAPLUS

CN Carbamic acid, [3-[(R)-[4-(cyclobutylmethyl)-1-piperazinyl][4-[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, methyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859843-98-4 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-methoxyethyl)-1-piperazinyl]methyl]phenyl]-, ethyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

● HC1

RN 859843-99-5 CAPLUS

CN Carbamic acid, [3-[(R)-[4-(cyclopropylmethyl)-1-piperazinyl][4-[(diethylamino)carbonyl]phenyl]methyl]phenyl]-, ethyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 859844-00-1 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl](4-propyl-1-piperazinyl)methyl]phenyl]-, ethyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

● HCl

RN 859844-01-2 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl](4-methyl-1-piperazinyl)methyl]phenyl]-, ethyl ester, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

IT 859635-21-5P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of diarylmethylpiperazines as $\boldsymbol{\delta}$ receptor ligands for treatment of pain)

RN 859635-21-5 CAPLUS

CN 1-Piperazinecarboxylic acid, 4-[(S)-[4-[(diethylamino)carbonyl]phenyl][3-[(methoxycarbonyl)amino]phenyl]methyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)

Absolute stereochemistry.

$$\begin{array}{c|c} & H \\ & N \\ & O \\ & &$$

REFERENCE COUNT: 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

ANSWER 2 OF 2 CAPLUS COPYRIGHT 2008 ACS on STN L3

ACCESSION NUMBER: 2004:412932 CAPLUS

DOCUMENT NUMBER: 140:423709

TITLE: Preparation of N-[4-(phenylpiperazinylmethyl)phenyl]ca

rbamates for treatment of pain, anxiety, or

gastrointestinal disorders

Brown, William; Griffin, Andrew; Jones, Paul; Page, INVENTOR(S):

Daniel; Plobeck, Niklas; Walpole, Christopher

PATENT ASSIGNEE(S): Astrazeneca AB, Swed. SOURCE: PCT Int. Appl., 72 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent English LANGUAGE:

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PA:	PATENT NO.						KIND DATE			APPLICATION NO.						DATE			
WO 2004041802																			
	W:	ΑE,	AG,	AL,	AM,	AT,	AU,	ΑZ,	BA,	BB,	BG,	BR,	BW,	BY,	BZ,	CA,	CH,		
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,		
		GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KP,	KR,	KZ,	LC,		
		LK,	LR,	LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	NI,	NO,		
		NZ,	OM,	PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,	SG,	SK,	SL,	SY,	ΤJ,		
		TM,	TN,	TR,	TT,	TZ	UA,	UG,	US,	UΖ,	VC,	VN,	YU,	ZA,	ZM,	ZW			
	RW:	BW,	GH,	GM,	KE,	LS,	MW,	MZ,	SD,	SL,	SZ,	TZ,	UG,	ZM,	ZW,	AM,	ΑZ,		
							TJ,												
		ES,	FI,	FR,	GB,	GR,	HU,	ΙE,	ΙΤ,	LU,	MC,	NL,	PT,	RO,	SE,	SI,	SK,		
							CI,											ΤG	
CA	2502	732			A1		2004	0521		CA 2003-2502732				20031105					
			A1 20040607				AU 2003-278665					20031105							
EP								EP 2003-770198											
	R:						ES,										PT,		
							RO,												
BR	BR 2003015995				A		2005	0927	BR 2003-15995					20031105					
	CN 1711252						2005	1221		CN 2	003-	8010	2831	20031105					
JP	JP 2006514002						2006	0427		.TP 2	0.04 -	5497	76	20031105					
NZ	NZ 539484 IN 2005DN01579					A 20070531				NZ 2	003-	5394	84	20031105					
IN	IN 2005DN01579						2006	1229		IN 2	005-	DN15	79	20050419					
	MX 2005PA04708																		
US	US 2006122193						2006	0608	US 2005-533654					20050504					
US	7253	173			В2		2007	0807											
ZA	5 7253173 A 2005003556				A 20060830			ZA 2005-3556											
	NO 2005002698																		
	US 2007254890					A1 20071101			US 2007-774935 SE 2002-3303										
PRIORIT	IORITY APPLN. INFO.:									SE 2	002-	3303		1	A 2	0021	107		
													07						
										US 2	005-	5336	54	1	A1 2	0050	504		
THER S	HER SOURCE(S):				MARPAT 140:42370				09										

GΙ

Title compds. I [wherein R1 = (un)substituted (hetero)aryl(alkyl); R2 and AΒ

^{*} STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

R3 = independently H or (un)substituted (cyclo)alkyl; or pharmaceutically acceptable salts, diastereomers, enantiomers, or mixts. thereof] were prepared as opioid δ receptor ligands. For example, 4-carboxybenzaldehyde was amidated with diethylamine using SOC12 in CH2C12 to give N, N-diethyl-4-formylbenzamide (90%). Coupling of the amide with N-Boc-piperazine in the presence of benzotriazole in toluene, followed by reaction with 3-bromophenylzinc iodide in THF, afforded tert-Bu 4-[(3-bromophenyl)[4-[(diethylamino)carbonyl]phenyl]methyl]-1piperazinecarboxylate (33%). Coupling with Me carbamate (62%) using xantphos, Cs2CO3, and Pd2(dba)3 in dioxane, deprotection (89%) with TFA in CH2Cl2, and chiral HPLC separation of the enantiomers provided (-)-[3-[[4-[(diethylamino)carbonyl]phenyl]-1-piperazinylmethyl]phenyl]carb amic acid Me ester. Reaction of the piperazine with benzaldehyde in the presence of NaBH(OAc)3 in CH2Cl2 gave (R)-II. In binding assays using human 293S cells expressing cloned human opioid receptors and neomycin resistance, (R)-II and ten other exemplified compds. exhibited strong binding to the δ receptor with IC50 values in the range of 0.25-0.74 nM and showed some activity toward the κ (IC50 = 247-1636 nM) and μ (IC50 = 93-1100 nM) receptors. In functional assays, (R)-II demonstrated δ receptor agonist activity by activating the binding of GTP to G-proteins. Thus, I and their pharmaceutical compns. are useful in therapy, in particular for the treatment of gastrointestinal disorders, anxiety, or pain (no data).

IT 691890-43-4P, 4-[[4-[(Diethylamino)carbonyl]phenyl][3[(methoxycarbonyl)amino]phenyl]methyl]-1-piperazinecarboxylic acid
1,1-dimethylethyl ester

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(intermediate; preparation of N-[(phenylpiperazinylmethyl)phenyl]carbamates as δ receptor agonists for treatment of pain, anxiety, or gastrointestinal disorders)

RN 691890-43-4 CAPLUS

CN

1-Piperazinecarboxylic acid, 4-[[4-[(diethylamino)carbonyl]phenyl][3-[(methoxycarbonyl)amino]phenyl]methyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)

IT 691890-67-2P

RL: PAC (Pharmacological activity); PUR (Purification or recovery); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(δ receptor agonist; preparation of N-[(phenylpiperazinylmethyl)phenyl

] carbamates as δ receptor agonists for treatment of pain, anxiety, or gastrointestinal disorders)

RN 691890-67-2 CAPLUS

CN Carbamic acid, [3-[[4-[(diethylamino)carbonyl]phenyl][4-(phenylmethyl)-1-piperazinyl]methyl]phenyl]methyl-, methyl ester, (-)- (9CI) (CA INDEX NAME)

Rotation (-).

IT 691890-72-9P 691890-74-1P 691890-76-3P
 RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
 (δ receptor agonist; preparation of N-[(phenylpiperazinylmethyl)phenyl]carbamates as δ receptor agonists for treatment of pain,
 anxiety, or gastrointestinal disorders)
RN 691890-72-9 CAPLUS
CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(4-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester,
 trifluoroacetate (5:9) (9CI) (CA INDEX NAME)

CRN 691890-71-8 CMF C28 H35 N5 O3 S

Absolute stereochemistry. Rotation (+).

CM 2

CRN 76-05-1 CMF C2 H F3 O2

RN 691890-74-1 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(5-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (5:13) (9CI) (CA INDEX NAME)

CM I

CRN 691890-73-0 CMF C28 H35 N5 O3 S

Absolute stereochemistry. Rotation (-).

CM 2

CRN 76-05-1 CMF C2 H F3 O2

RN 691890-76-3 CAPLUS
CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(5-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (9CI) (CA INDEX NAME)

CM I

CRN 691890-75-2 CMF C28 H35 N5 O3 S

Absolute stereochemistry. Rotation (+).

CM 2

CRN 76-05-1 CMF C2 H F3 O2

```
691890-51-4P, (R)-Methyl [3-[[4-[(diethylamino)carbonyl]phenyl](4-
ΤТ
    benzylpiperazin-1-yl)methyl]phenyl]carbamate 691890-52-5P,
     (S)-Methyl [3-[[4-[(diethylamino)carbonyl]phenyl](4-benzylpiperazin-1-
    yl)methyl]phenyl]carbamate 691890-53-6P 691890-54-7P
    691890-55-8P 691890-56-9P 691890-57-0P
    691890-58-1P 691890-59-2P 691890-60-5P
    691890-61-6P 691890-62-7P 691890-63-8P
    691890-64-9P 691890-65-0P 691890-66-1P
    691890-68-3P 691890-69-4P 691890-70-7P
    691890-71-8P 691890-73-0P 691890-75-2P
    691890-77-4P, Methyl [3-[[4-[(diethylamino)carbonyl]phenyl](4-
    benzylpiperazin-1-yl)methyl]phenyl]carbamate 691890-78-5P,
    Methyl [3-[[4-[(diethylamino)carbonyl]phenyl][4-(thien-2-
    ylmethyl)piperazin-1-yl]methyl]phenyl]carbamate 691890-79-6P,
    Methyl [3-[[4-[(diethylamino)carbonyl]phenyl][4-(thien-3-
    ylmethyl)piperazin-1-yl]methyl]phenyl]carbamate 691890-80-9P,
    Methyl [3-[[4-[(diethylamino)carbonyl]phenyl][4-(2-furylmethyl)piperazin-1-
    yl]methyl]phenyl]carbamate 691890-81-0P, Methyl
    [3-[4-[diethylamino)carbonyl]phenyl][4-(3-furylmethyl)piperazin-1-
    yl]methyl]phenyl]carbamate 691890-82-1P, Methyl
    [3-[[4-[(diethylamino)carbonyl]phenyl][4-(1H-imidazol-2-ylmethyl)piperazin-
    1-y1]methy1]pheny1]carbamate 691890-83-2P, Methy1
    [3-[[4-[(diethylamino)carbonyl]phenyl]][4-(pyridin-2-ylmethyl)piperazin-1-
    yl]methyl]phenyl]carbamate 691890-84-3P, Methyl
    [3-[[4-[(diethylamino)carbonyl]phenyl][4-(pyridin-4-ylmethyl)piperazin-1-
    yl]methyl]phenyl]carbamate 691890-85-4P, Methyl
```

```
[3-[[4-[(diethylamino)carbonyl]phenyl][4-(1,3-thiazol-2-ylmethyl)piperazin-
     1-y1]methy1]pheny1]carbamate 691890-86-5P 691890-87-6P
     691890-88-7P 691890-89-8P 691890-90-1P
     691890-91-2P 691890-92-3P 691890-93-4P
     691890-94-5P 691890-95-6P 691890-96-7P
     691890-97-8P 691890-98-9P 691890-99-0P
     691891-00-6P 691891-01-7P 691891-02-8P
     691891-03-9P
     RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
     (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
     (Uses)
        (\delta receptor agonist; preparation of N-[(phenylpiperazinylmethyl)phenyl
        ] carbamates as \delta receptor agonists for treatment of pain,
        anxiety, or gastrointestinal disorders)
     691890-51-4 CAPLUS
RN
     Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(phenylmethyl)-
CN
     1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)
```

Absolute stereochemistry. Rotation (-).

RN 691890-52-5 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(phenylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-53-6 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(2-thienylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-54-7 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-thienylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 691890-55-8 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(3-thienylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-56-9 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(3-thienylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 691890-57-0 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(2-furanylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-58-1 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-furanylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 691890-59-2 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(3furanylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-60-5 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(3-furanylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 691890-61-6 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(1H-imidazol-2-ylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 691890-62-7 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(2-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-63-8 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 691890-64-9 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(4-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-65-0 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(4-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 691890-66-1 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 691890-68-3 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(3-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-69-4 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(2-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-70-7 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(4-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 691890-71-8 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(4-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-73-0 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(5-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 691890-75-2 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(5-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 691890-77-4 CAPLUS

CN Carbamic acid, [3-[[4-[(diethylamino)carbonyl]phenyl][4-(phenylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

RN 691890-78-5 CAPLUS

CN Carbamic acid, [3-[[4-[(diethylamino)carbonyl]phenyl][4-(2-thienylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

RN 691890-79-6 CAPLUS

CN Carbamic acid, [3-[[4-[(diethylamino)carbonyl]phenyl][4-(3-thienylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

RN 691890-80-9 CAPLUS

CN Carbamic acid, [3-[[4-[(diethylamino)carbonyl]phenyl][4-(2-furanylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

RN 691890-81-0 CAPLUS

CN Carbamic acid, [3-[[4-[(diethylamino)carbonyl]phenyl][4-(3-furanylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

RN 691890-82-1 CAPLUS

CN Carbamic acid, [3-[[4-[(diethylamino)carbonyl]phenyl][4-(1H-imidazol-2-ylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

RN 691890-83-2 CAPLUS

CN Carbamic acid, [3-[[4-[(diethylamino)carbonyl]phenyl]][4-(2-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

RN 691890-84-3 CAPLUS

CN Carbamic acid, [3-[[4-[(diethylamino)carbonyl]phenyl]][4-(4-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CA INDEX NAME)

RN 691890-85-4 CAPLUS

CN Carbamic acid, [3-[[4-[(diethylamino)carbonyl]phenyl][4-(2-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester (9CI) (CAINDEX NAME)

RN 691890-86-5 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(phenylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (10:21) (9CI) (CA INDEX NAME)

CRN 691890-51-4 CMF C31 H38 N4 O3

Absolute stereochemistry. Rotation (-).

CM 2

CRN 76-05-1 CMF C2 H F3 O2

RN 691890-87-6 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(phenylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (5:8) (9CI) (CA INDEX NAME)

CM 1

CRN 691890-52-5 CMF C31 H38 N4 O3

Absolute stereochemistry. Rotation (+).

<12/04/2007>

Erich Leese

CRN 76-05-1 CMF C2 H F3 O2

$${\tiny \begin{array}{c} F\\ |\\ F-C-CO_2H\\ |\\ F\end{array}}$$

RN 691890-88-7 CAPLUS
CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(2-thienylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (5:7) (9CI) (CA INDEX NAME)

CM 1

CRN 691890-53-6 CMF C29 H36 N4 O3 S

Absolute stereochemistry. Rotation (+).

CRN 76-05-1 CMF C2 H F3 O2

RN 691890-89-8 CAPLUS
CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-thienylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (2:3) (9CI) (CA INDEX NAME)

CM I

CRN 691890-54-7 CMF C29 H36 N4 O3 S

Absolute stereochemistry. Rotation (-).

CRN 76-05-1 CMF C2 H F3 O2

$${\tiny \begin{array}{c}F\\F-C-CO_2H\\\\F\end{array}}$$

691890-90-1 CAPLUS RN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(3-CN

thienylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (10:17) (9CI) (CA INDEX NAME)

CM

CRN 691890-55-8

CMF C29 H36 N4 O3 S

Absolute stereochemistry. Rotation (+).

CRN 76-05-1 CMF C2 H F3 O2

$${\tiny \begin{array}{c}F\\ \\F-C-CO_2H\\ \\F\end{array}}$$

RN 691890-91-2 CAPLUS
CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(3-thienylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (5:7) (9CI) (CA INDEX NAME)

CM I

CRN 691890-56-9 CMF C29 H36 N4 O3 S

Absolute stereochemistry. Rotation (-).

CRN 76-05-1 CMF C2 H F3 O2

RN 691890-92-3 CAPLUS
CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(2-furanylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (2:3) (9CI) (CA INDEX NAME)

CM 1

CRN 691890-57-0 CMF C29 H36 N4 O4

Absolute stereochemistry. Rotation (+).

CRN 76-05-1 CMF C2 H F3 O2

RN 691890-93-4 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-furanylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (5:8) (9CI) (CA INDEX NAME)

CM 1

CRN 691890-58-1 CMF C29 H36 N4 O4

Absolute stereochemistry. Rotation (-).

CRN 76-05-1 CMF C2 H F3 O2

RN 691890-94-5 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(3-furanylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, bis(trifluoroacetate) (9CI) (CA INDEX NAME)

CM I

CRN 691890-59-2 CMF C29 H36 N4 O4

Absolute stereochemistry. Rotation (+).

CRN 76-05-1 CMF C2 H F3 O2

RN 691890-95-6 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(1H-imidazol-2-ylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (10:21) (9CI) (CA INDEX NAME)

CM 1

CRN 691890-61-6 CMF C28 H36 N6 O3

Absolute stereochemistry. Rotation (-).

CRN 76-05-1 CMF C2 H F3 O2

RN 691890-96-7 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(2-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (2:3) (9CI) (CA INDEX NAME)

CM 1

CRN 691890-62-7 CMF C30 H37 N5 O3

Absolute stereochemistry. Rotation (+).

CRN 76-05-1 CMF C2 H F3 O2

$${\tiny \begin{array}{c}F\\F-C-CO_2H\\\\F\end{array}}$$

RN 691890-97-8 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(2-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (10:17) (9CI) (CA INDEX NAME)

CM 1

CRN 691890-63-8 CMF C30 H37 N5 O3

Absolute stereochemistry. Rotation (-).

CRN 76-05-1 CMF C2 H F3 O2

RN 691890-98-9 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(4-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (10:19) (9CI) (CA INDEX NAME)

CM I

CRN 691890-64-9 CMF C30 H37 N5 O3

Absolute stereochemistry. Rotation (+).

CRN 76-05-1 CMF C2 H F3 O2

$${\tiny \begin{array}{c}F\\ -\\ C-CO_2H\\ -\\ F\end{array}}$$

RN 691890-99-0 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(4-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (5:13) (9CI) (CA INDEX NAME)

CM I

CRN 691890-65-0 CMF C30 H37 N5 O3

Absolute stereochemistry. Rotation (-).

CRN 76-05-1 CMF C2 H F3 O2

RN 691891-00-6 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(3-pyridinylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, hydrochloride (10:29) (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

●29/10 HC1

RN 691891-01-7 CAPLUS

CN Carbamic acid, [3-[(S)-[4-[(diethylamino)carbonyl]phenyl][4-(2-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, hydrochloride (2:3) (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

●3/2 HC1

RN 691891-02-8 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(4-thiazolylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (9CI) (CA INDEX NAME)

CM 1

CRN 691890-70-7 CMF C28 H35 N5 O3 S

Absolute stereochemistry. Rotation (-).

<12/04/2007>

10/513699

CM 2

CRN 76-05-1 CMF C2 H F3 O2

RN 691891-03-9 CAPLUS

CN Carbamic acid, [3-[(R)-[4-[(diethylamino)carbonyl]phenyl][4-(3-furanylmethyl)-1-piperazinyl]methyl]phenyl]-, methyl ester, trifluoroacetate (2:3) (9CI) (CA INDEX NAME)

CM 1

CRN 691890-60-5 CMF C29 H36 N4 O4

Absolute stereochemistry. Rotation (-).

$$\begin{array}{c|c} \text{Et}_2N & \text{O} \\ \hline \\ N & R \\ \hline \end{array}$$

CM 2

CRN 76-05-1 CMF C2 H F3 O2

=> d his

(FILE 'HOME' ENTERED AT 15:01:03 ON 18 JAN 2008)

FILE 'REGISTRY' ENTERED AT 15:01:15 ON 18 JAN 2008

L1 STRUCTURE UPLOADED

L2 90 S L1 FULL

FILE 'CAPLUS' ENTERED AT 15:02:19 ON 18 JAN 2008

L3 2 S L2 FULL

=> log y

COST IN U.S. DOLLARS

SINCE FILE
ENTRY
SESSION
FULL ESTIMATED COST

12.82

TOTAL
191.85

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE -1.60 -1.60

STN INTERNATIONAL LOGOFF AT 15:04:33 ON 18 JAN 2008